63377 SEARCH REQUEST FORM

	Requestor's Name:	akslimi C	bannalajai	Serial — Number: (09/774,55	8
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=> fil reg FILE 'REGISTRY' ENTERED AT 07:10:08 ON 29 MAR 2002 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2002 American Chemical Society (ACS)

STRUCTURE FILE UPDATES: 27 MAR 2002 HIGHEST RN 403476-73-3 DICTIONARY FILE UPDATES: 27 MAR 2002 HIGHEST RN 403476-73-3

TSCA INFORMATION NOW CURRENT THROUGH July 7, 2001

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Calculated physical property data is now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details: http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf

The P indicator for Preparations was not generated for all of the CAS Registry Numbers that were added to the H/Z/CA/CAplus files between 12/27/01 and 1/23/02. Use of the P indicator in online and SDI searches during this period, either directly appended to a CAS Registry Number or by qualifying an L-number with /P, may have yielded incomplete results. As of 1/23/02, the situation has been resolved. Also, note that searches conducted using the PREP role indicator were not affected.

Customers running searches and/or SDIs in the H/Z/CA/CAplus files incorporating CAS Registry Numbers with the P indicator between 12/27/01 and 1/23/02, are encouraged to re-run these strategies. Contact the CAS Help Desk at 1-800-848-6533 in North America or 1-614-447-3698, worldwide, or send an e-mail to help@cas.org for further assistance or to receive a credit for any duplicate searches.

VAR G1=O/21 VAR G4=O/S VAR G5=O/S/N NODE ATTRIBUTES: DEFAULT MLEVEL IS ATOM DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES: RING(S) ARE ISOLATED OR EMBEDDED NUMBER OF NODES IS 18

STEREO ATTRIBUTES: NONE

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Jan Delaval Reference Librarian Biotechnology & Chemical Library CM1 1E07 – 703-308-4498 jan.delaval@uspto.gov

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100.0% PROCESSED 92 ITERATIONS SEARCH TIME: 00.00.02
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90 ANSWERS

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     (FILE 'HCAPLUS' ENTERED AT 06:38:21 ON 29 MAR 2002)
                DEL HIS
                E WO2001-US10369/AP, PRN
              1 S E3
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          10544 S (PROC?(L)GAMB?)/PA,CS
L2
                E DELONG M/AU
             43 S E3, E4, E21-E25
L3
                E DE LONG M/AU
              5 S E7
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                E MCIVER J/AU
             54 S E3, E6, E16, E18, E19
L5
                E MC IVER J/AU
                E MCMILLAN J/AU
             30 S E3, E14, E42
L6
                E MC MILLAN J/AU
                E YOUNGQUIST R/AU
             82 S E3, E5, E6, E9, E11, E12
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                SEL RN L1
     FILE 'REGISTRY' ENTERED AT 06:47:58 ON 29 MAR 2002
L11
             51 S E1-E51
             43 S L11 AND 16.127/RID
L12
              8 S L11 NOT L12
L13
                E 16.127/RID
           7268 S E3 AND P/ELS
L14
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                STR L9
              4 S L15 SAM SUB=L14
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L21
             43 S L11 AND L20
L22
             47 S L20 NOT L21
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L24
             34 S L22 NOT L23
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L28
              1 S L21
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             15 S L24
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              3 S L30 AND L28, L29
L32
              6 S L28, L29, L31
L33
              4 S L32 AND L1-L7
L34
              6 S L32, L33
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FILE 'USPATFULL, USPAT2' ENTERED AT 07:09:17 ON 29 MAR 2002

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L35
              1 S L21
L36
              1 S L23
              8 S L24
L37
L38
              0 S L37 AND L35, L36
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              2 S L35, L36
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              7 DUP REM L34 L39 (1 DUPLICATE REMOVED)
     FILE 'REGISTRY' ENTERED AT 07:10:08 ON 29 MAR 2002
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FILE 'HCAPLUS' ENTERED AT 07:10:30 ON 29 MAR 2002
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PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
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FILE 'USPATFULL' ENTERED AT 07:10:30 ON 29 MAR 2002
CA INDEXING COPYRIGHT (C) 2002 AMERICAN CHEMICAL SOCIETY (ACS)
FILE 'USPAT2' ENTERED AT 07:10:30 ON 29 MAR 2002
CA INDEXING COPYRIGHT (C) 2002 AMERICAN CHEMICAL SOCIETY (ACS)
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    ANSWER 1 OF 7 HCAPLUS COPYRIGHT 2002 ACS
                                                        DUPLICATE 1
L40
     1980:58302 HCAPLUS
AN
     92:58302
DN
TI
     1 and 2-Substituted analogs of certain prostaglandins
IN
     Biddlecom, William G.; Kluender, Harold C.; Woessner, Warren D.
PΑ
     Miles Laboratories, Inc., USA
     U.S., 40 pp.
SO
     CODEN: USXXAM
DT
     Patent
     English
LA
FAN.CNT 1
                     KIND DATE
                                          APPLICATION NO. DATE
     PATENT NO.
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* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

US 1978-912515

19780605

A

19791016

- AB App. 110 decarboxyprostaglandin analogs and intermediates for them were prepd. from the key compds. I (n=6,7) by appropriate modifications of conventional syntheses; extensive biol. activity tests (inhibition of platelet aggregation, and gastric secretion, antihyprotensive and muscle relaxant activity) were carried out. Among the compds. prepd. were II-V.
- TT 72488-21-2P 72489-45-3P
 RL: BAC (Biological activity or effector, except adverse); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation)
 (prepn. and biol. activity of)
- CN Phosphonic acid, [(9.alpha.,11.alpha.,13E,15S)-9,11,15-trihydroxyprost-13-en-1-yl]-, dimethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Double bond geometry as shown.

PΙ

GΙ

US 4171331

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ANSWER 2 OF 7 HCAPLUS COPYRIGHT 2002 ACS
L40
ΑN
     2001:747564 HCAPLUS
     135:293970
DN
ΤI
     Cosmetic and pharmaceutical compositions and methods using
     2-decarboxy-2-phosphinico prostaglandin derivatives
                                                                    references
     Delong, Mitchell Anthony; Mciver, John Mcmillan;
IN
     Youngquist, Robert Scott
     The Procter + Gamble Company, USA
PA
     PCT Int. Appl., 54 pp.
SO
     CODEN: PIXXD2
DT
     Patent
     English
LA
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     PATENT NO.
                      KIND
                            DATE
                                           APPLICATION NO.
                                                            DATE
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                            20011011
                                           WO 2001-US10369
     WO 2001074314
ΡI
                            20020221
     WO 2001074314
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            AE, AG, AL, AM, AT, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH,
             CN, CO, CR, CU, CZ, CZ, DE, DE, DK, DK, DM, DZ, EE, EE, ES, FI,
             FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP,
             KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX,
             MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SK, SL, TJ, TM,
             TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD,
             RU, TJ, TM
         RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY,
             DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF,
             BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG
                                           US 2001-774558
                                                            20010131
     US 2002013294
                       Α1
                            20020131
PRAI US 2000-193845P
                            20000331
     MARPAT 135:293970
OS
     Compns. contq. 2-decarboxy-2-phosphinico prostaglandin derivs. is
AB
     described for treating hair loss in mammals. The compns. can be applied
     topically to the skin to arrest hair loss, reverse hair loss, and promote
     hair growth. Compns. contg. 2-decarboxy-2-phosphinico prostaglandin
     derivs. can also be used to lower intraocular pressure and treat bone
     disorders. A compn. comprises a prostaglandin analog, an activity
     enhancer, such as a hair growth stimulant and a penetration enhancer, and
     a sufficient amt. of a component selected from the group consisting of
     emollients, propellants, solvents, humectants, thickeners, powders,
     fragrances, water, alcs., aloe vera gel, allantoin, glycerin, vitamin A
     and E oils, mineral oil, propylene glycol, polypropylene glycol-2 myristyl
     propionate, di-Me isosorbide, and combinations thereof. For example, a
     compn. for topical administration was prepd. comprising (by wt.) a
     prostaglandin (IC50 = 114 nM) 1.14%, ethanol 59.32%, propylene glycol
     19.77%, and di-Me isosorbide 19.77%. Also, a shampoo was made contg.
     ammonium lauryl sulfate 11.5%, ammonium laureth sulfate 4%, cocamide MEA
     2%, ethylene glycol distearate 2%, cetyl alc. 2%, stearyl alc. 1.2%,
     glycerin 1%, sodium chloride 0.1%, sucrose polyesters of cottonate fatty
     acid 3%, sucrose polyesters of behenate fatty acid 2%, lauryl di-Me amine
     oxide 1.5%, DMDM hydantoin 0.15%, prostaglandin (IC = 150 nM) 0.15%,
     phenoxyethanol 0.5%, fragrance 0.5%, and water up to 100%. A tablet
```

formulation was also prepd. contg. a prostaglandin 5 mg, microcryst. cellulose 100 mg, sodium starch glycolate 30 mg, and magnesium stearate 3 mg per tablet. When administered orally once daily, the above compn. substantially increases bone vol. in a patient suffering from osteoporosis.

TT 365241-18-5P 365241-19-6P 365241-20-9P 365241-21-0P 365241-22-1P 365241-23-2P 365241-24-3P 365241-25-4P 365241-26-5P 365241-27-6P 365241-28-7P 365241-29-8P 365241-30-1P 365241-31-2P 365241-32-3P 365241-33-4P 365241-34-5P 365241-35-6P 365241-36-7P 365241-37-8P 365241-38-9P 365241-42-5P 365241-40-3P 365241-41-4P 365241-45-8P 365241-43-6P 365241-47-0P 365241-48-1P 365241-49-2P 365241-50-5P 365241-51-6P 365241-55-0P 365241-56-1P 365241-57-2P 365241-58-3P 365241-59-4P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); BUU (Biological use, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(cosmetic and pharmaceutical compns. contg. 2-decarboxy-2-phosphinico prostaglandin derivs.)

IT 365241-18-5P

365241-60-7P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); BUU (Biological use, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(cosmetic and pharmaceutical compns. contg. 2-decarboxy-2-phosphinico prostaglandin derivs.)

RN 365241-18-5 HCAPLUS

CN Phosphinic acid, [6-[(1R,5S)-2-[(3R)-4-[(3-fluorophenyl)thio]-3-hydroxybutyl]-5-hydroxy-3-(hydroxyimino)cyclopentyl]hexyl]methyl- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Double bond geometry unknown.

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ANSWER 3 OF 7 HCAPLUS COPYRIGHT 2002 ACS
L40
ΑN
     2001:115155 HCAPLUS
DN
     134:162868
     Novel 2-decarboxy-2-phosphinico prostaglandin F analogs
ΤI
     Delong, Mitchell Anthony; Wos, John August; De, Biswanath;
IN
     Ebetino, Frank Hallock
PA
     Procter & Gamble Company, USA
SO
     PCT Int. Appl., 35 pp.
     CODEN: PIXXD2
DT
     Patent
LA
     English
FAN.CNT 1
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			CN,	CR,	CU,	CZ,	CZ,	DE,	DE,	DK,	DK,	DM,	DZ,	EE,	EE,	ES,	FI,	FI,
			GB,	GD,	GE,	GH,	GM,	HR,	HU,	ID,	IL,	IN,	IS,	JP,	KΕ,	KG,	KΡ,	KR,
															MK,			
			ΜZ,	NO,	NZ,	PL,	PT,	RO,	RU,	SD,	SE,	SG,	SI,	SK,	SK,	SL,	ТJ,	TM,
			TR,	TT,	ΤZ,	UA,	UG,	UZ,	VN,	YU,	ZA,	ZW,	AM,	ΑZ,	BY,	KG,	ΚZ,	MD,
				ΤJ,														
		RW:	GH,	GM,	KΕ,	LS,	MW,	ΜZ,	SD,	SL,	SZ,	TZ,	UG,	ZW,	AT,	ΒE,	CH,	CY,
			•				-								PT,	SE,	BF,	ВJ,
							GA,		GW,	\mathtt{ML} ,	MR,	ΝE,	SN,	TD,	TG			
PRAI	US	1999	-147	132P	P		1999	0804										
os	MAI	RPAT :	134:	1628	68													
GI																		

search of

HO
$$X$$
 HO $Y - R^1$ $Y - R^2$ HO $Y - R^2$

AB Novel prostaglandin F analogs of formula I [R1 = H, alkyl; R2 = H, alkyl, heteroalkyl, carbocyclic ring; X = O, S; Y = O, S, NH; V = alkyl, heteroalkyl; W = OH, (substituted) NHOH; (substituted) NOH; Z = alkyl, heteroalkyl, alkyl-cycloalkyl, etc.] are prepd. This invention also includes optical isomers, diastereomers and enantiomers of the formula, and pharmaceutically-acceptable salts, biohydrolyzable amides, esters, and imides thereof. The compds. of the present invention are useful for the treatment of a variety of diseases and conditions, such as bone disorders and glaucoma. Accordingly, the invention further provides pharmaceutical compns. comprising these compds. The invention still further provides methods of treatment for bone disorders and glaucoma using these compds. or the compns. contg. them. Thus, II is prepd. and is used in tablets and liq. compns.

IT 325153-11-5P 325153-12-6P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(prepn. of prostaglandin F analogs for the treatment of osteoporosis and glaucoma) $\,$

IT 325153-11-5P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(prepn. of prostaglandin F analogs for the treatment of osteoporosis and glaucoma)

RN 325153-11-5 HCAPLUS

CN Phosphinic acid, [(4Z)-6-[(1R,2R,3R,5S)-2-[(1E,3S)-5-(2-fluorophenyl)-3-hydroxy-1-pentenyl]-3,5-dihydroxycyclopentyl]-4-hexenyl]methyl-, monosodium salt (9CI) (CA INDEX NAME)

Absolute stereochemistry.
Double bond geometry as shown.

Na

RETABLE Referenced Author Year (RAU) (RPY) (RVL) (RPG)	(RWK)	Referenced File
Kende, A 1999 Kluender, H 1979 Liljebris, C 1995 The Procter & Gamble Co 1999 The Procter & Gamble Co 1999 The Procter & Gamble Co 1999	40 8189 2 441 38 289 	TETRAHEDRON LETTERS PROSTAGLANDINS MED JOURNAL OF MEDICINAL WO 9912895 A WO 9912896 A WO 9912898 A	HCAPLUS HCAPLUS
L40 ANSWER 4 OF 7 HCAPLUS AN 2001:716927 HCAPLUS	COPYRIGHT 200	02 ACS	

DN 136:31838

Synthesis and biological evaluation of prostaglandin-F alkylphosphinic ΤI acid derivatives as bone anabolic agents for the treatment of osteoporosis ΑU

Soper, David L.; Milbank, Jared B. J.; Mieling, Glen E.; Dirr, Michelle J.; Kende, Andrew S.; Cooper, Robin; Jee, Webster S. S.; Yao, Wei; Chen, Jian Liang; Bodman, Mark; Lundy, Mark W.; De, Biswanath; Stella, Mark E.; Ebetino, Frank H.; Wang, Yili; de Long, Mitchell A.; Wos, John

X

Procter & Gamble Pharmaceuticals, Mason, OH, 45040, USA CS SO

Journal of Medicinal Chemistry (2001), 44(24), 4157-4169

CODEN: JMCMAR; ISSN: 0022-2623

PB American Chemical Society

DTJournal

LA English

GΙ

A series of novel C1 alkylphosphinic acid analogs of the prostaglandin-F AΒ family have been evaluated at the eight human prostaglandin receptors for potential use in the treatment of osteoporosis. Using mol. modeling as a tool for structure-based drug design, the authors have discovered that the phosphinic acid moiety (P(O)(OH)R) behaves as an isostere for the C1

Ι

carboxylic acid in the human prostaglandin FP binding assay in vitro and possesses enhanced hFP receptor selectivity when compared to the parent carboxylic acid. When evaluated in vivo, the Me phosphinic acid analog (I) produced a bone anabolic response in rats, returning bone mineral d. (BMD) to intact levels in the distal femur in the ovariectomized rat (OVX) model. These results suggest that prostaglandins of this class may be useful agents in the treatment of diseases assocd. with bone loss. 380153-29-7P

RL: DMA (Drug mechanism of action); PAC (Pharmacological activity); PRP (Properties); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)

(synthesis and biol. evaluation of prostaglandin F alkylphosphinic acid derivs. as bone anabolic agents for treatment of osteoporosis)

IT 252858-16-5P 380153-33-3P 380153-34-4P 380153-35-5P 380153-36-6P 380153-37-7P

RL: DMA (Drug mechanism of action); PAC (Pharmacological activity); PRP (Properties); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(synthesis and biol. evaluation of prostaglandin F alkylphosphinic acid derivs. as bone anabolic agents for treatment of osteoporosis)

IT 380153-32-2P 380153-41-3P 380153-42-4P 380153-43-5P 380153-44-6P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(synthesis and biol. evaluation of prostaglandin F alkylphosphinic acid derivs. as bone anabolic agents for treatment of osteoporosis)

IT 380153-29-7P

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RN

RL: DMA (Drug mechanism of action); PAC (Pharmacological activity); PRP (Properties); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)

(synthesis and biol. evaluation of prostaglandin F alkylphosphinic acid derivs. as bone anabolic agents for treatment of osteoporosis)

380153-29-7 HCAPLUS

CN Phosphinic acid, [(4Z)-6-[(1R,2R,3R,5S)-3,5-dihydroxy-2-[(1E,3R)-3-hydroxy-4-phenoxy-1-butenyl]cyclopentyl]-4-hexenyl]methyl- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Double bond geometry as shown.

RETABLE |Year | VOL | PG | Referenced Work | Referenced Referenced Author | (RPY) | (RVL) | (RPG) | (RWK) | File ________________________________ |Using ACD/pKa DB S| Advanced Chemistry Devel IHCAPLUS |1985 |41 3901 |Tetrahedron Barton, D |1983 |24 |4979 |Tetrahedron Lett IHCAPLUS Barton, D |2001 |265 1101 |Anat Rec (New Anat) | MEDLINE Borah, B | J Chem Soc, Chem Com | HCAPLUS |1992 | 1355 Bricklebank, N |521 |Drugs Future Cascierci, M |1996 |21 |1268 |J Clin Invest IMEDLINE Civitelli, R |1988 |82

Corey, É	1981		3455	Tetrahedron Lett	
Cummings, S	1990	•	163	Clin Orthop Relat Re	
deLong, M	12000		1519	Bioorg Med Chem Lett	
Engel, R	1988	1	21	Synthesis of Carbon-	!
Ericksen, E		16	381	Bone	<u> </u>
Francis, R	•	156	831	Curr Ther Res Clin E	
Funk, C	•	4 4	934		HCAPLUS
Furet, P		10	2337	Bioorgan Med Chem Le	
Gowen, M		15	1		HCAPLUS
GraphPad Software Inc	1999	1	1	The ED50 and 95% con	
Jee, W	1997	21	1297	• =	HCAPLUS
Jee, W			1101	Handbook of Bone Mor	
Ke, H	11998	123	1249	Bone	HCAPLUS
Kende, A	11999	40	18189	Tetrahedron Lett	HCAPLUS
Liljebris, C	11995	138	1289	J Med Chem	HCAPLUS
Liljebris, C	1996	161	14028	J Org Chem	HCAPLUS
Lindsay, R	11993	177	1535	J Clin Endocrinol Me	MEDLINE
Lopez, F	12000	4	1383	Curr Opin Chem Biol	HCAPLUS
Ma, Y	11995	17	549	Bone	HCAPLUS
Machwate, M	2001	60	36	Mol Pharmacol	HCAPLUS
Melton, L	11995	110	175	J Bone Miner Res	l
Mitlak, B	11997	148	155	Horm Res	HCAPLUS
Mohan, S	12000	27	471	Bone	HCAPLUS
Parfitt, A	1987	2	595	J Bone Miner Res	MEDLINE
Parfitt, A		172	1396	J Clin Invest	MEDLINE
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Sandri, J	11992	122	2945	Synth Commun	HCAPLUS
Schaaf, T	1979	122	1340	J Med Chem	HCAPLUS
Schaaf, T	11981	124	1353	J Med Chem	HCAPLUS
Stjernschantz, J	1992	17	691	Drugs Future	1
Turner, R	•	15	1275		HCAPLUS
Ueda, K		97	1834	J Pediatr	MEDLINE
Ueno, K		16	179	Bone	HCAPLUS
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Wang, Y	2000	•	1332		HCAPLUS
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L40 ANSWER 5 OF 7 HCAPLUS COPYRIGHT 2002 ACS

AN 1999:719514 HCAPLUS

DN 132:49813

Prostaglandin phosphonic acids through homolytic halodecarboxylation of TI prostaglandins F1.alpha. and F2.alpha.

Kende, Andrew S.; Milbank, Jared B. J.; Ebetino, Frank H.; Delong, AU Mitchell A.

Department of Chemistry, University of Rochester, Rochester, NY, 14627, CS

Tetrahedron Lett. ((1999)), 40(47), 8189-8192 SO CODEN: TELEAY; ISSN: 0040-4039

PΒ Elsevier Science Ltd.

DTJournal LA English

OS CASREACT 132:49813

GΙ

Published 1999 19 Nov 1999

AB Phosphonic acid derivs., e.g. I and II, of prostaglandins F1.alpha. and F2.alpha. were prepd. through Arbuzov reaction of 2-decarboxy-2-iodoprostaglandin intermediates. The intermediate iodo compds., which are potentially valuable for the synthesis of other analogs, were obtained from the parent prostaglandins by Barton's modification of the Hunsdiecker reaction.

IT 252858-11-0P 252858-14-3P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation) (prepn. of prostaglandin phosphonic acids through homolytic halodecarboxylation of prostaglandins F1.alpha. and F2.alpha.)

IT 252858-12-1P 252858-16-5P 252858-17-6P

RL: SPN (Synthetic preparation); PREP (Preparation) (prepn. of prostaglandin phosphonic acids through homolytic halodecarboxylation of prostaglandins F1.alpha. and F2.alpha.)

IT 252858-11-0P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation) (prepn. of prostaglandin phosphonic acids through homolytic halodecarboxylation of prostaglandins F1.alpha. and F2.alpha.)

RN 252858-11-0 HCAPLUS

CN Phosphonic acid, [6-[(1R,2R,3R,5S)-3,5-bis(acetyloxy)-2-[(1E,3S)-3-(acetyloxy)-1-octenyl]cyclopentyl]hexyl]-, diethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.
Double bond geometry as shown.

RETABLE

Referenced Author (RAU)	Year VOL (RPY) (RVL)	(RPG)	Referenced Work Referenced (RWK) File
	=+=====	+=	-+
Anon	1982	1	New Synthetic Routes

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|1985 |41
                                       3901
                                              |Tetrahedron
                                                                      | HCAPLUS
Barton, D
Barton, D
                         |1983 |24
                                       14979
                                              |Tetrahedron Lett
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                         |1977 |
                                              |Prostaglandin Synthe|
Bindra, J
Collins, P
                         11993 | 93
                                       11533
                                              |Chem Rev
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                         |1988 |
                                       |21
                                              |Synthesis of Carbon-|
Engel, R
                         |1975 |40
                                       |521
                                              | J Org Chem
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Iguchi, Y
                                       1219
                                              |Chem Rev
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Johnson, R
                         |1956 |56
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Kluender, H
                         |1973 |38
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                                                                      | HCAPLUS
Lincoln, F
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                                              |Tetrahedron Lett
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McKenna, C
Tsanaktidis, J
                         |1989 |30
                                       |6967
                                              |Tetrahedron Lett
                                       |332
                                              |Org React
                         |1957 | 9
Wilson, C
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L40 ANSWER 6 OF 7 HCAPLUS COPYRIGHT 2002 ACS

AN 1981:191740 HCAPLUS

DN 94:191740

TI The synthesis of dimethylphosphonoprostaglandin analogs

AU Kluender, Harold C.; Woessner, Warren

CS Miles Lab., Inc., Madison, WI, 53704, USA

SO Prostaglandins Med. (1979), 2(6), 441-4 CODEN: PROMDL; ISSN: 0161-4630

DT Journal

LA English

AB Several known title compds. were prepd. conventionally.

IT 77493-13-1P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation) (prepn. and deprotection of)

IT 72488-21-2P 72489-45-3P

RL: SPN (Synthetic preparation); PREP (Preparation)

(prepn. of)

IT 77493-13-1P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation) (prepn. and deprotection of)

RN 77493-13-1 HCAPLUS

CN Phosphonic acid, [(9.alpha.,11.alpha.,13E,15S)-9,11,15-tris[(tetrahydro-2H-pyran-2-yl)oxy]prost-13-en-1-yl]-, dimethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.
Double bond geometry as shown.

L40 ANSWER 7 OF 7 USPATFULL

AN 2002:22458 USPATFULL

TI Cosmetic and pharmaceutical compositions and methods using

2-decarboxy-2-phosphinico derivatives

IN DeLong, Mitchell Anthony, West Chester, OH, UNITED STATES McIver, John McMillan, Cincinnati, OH, UNITED STATES

Youngquist, Robert Scott, Mason, OH, UNITED STATES

PI US 2002013294 A1 20020131 AI US 2001-774558 A1 20010131 (9) PRAI US 2000-193845P 20000331 (60)

DT Utility
FS APPLICATION

LREP THE PROCTER & GAMBLE COMPANY, PATENT DIVISION, IVORYDALE TECHNICAL CENTER - BOX 474, 5299 SPRING GROVE AVENUE, CINCINNATI, OH, 45217

CLMN Number of Claims: 30 ECL Exemplary Claim: 1

DRWN No Drawings

LN.CNT 1847

CAS INDEXING IS AVAILABLE FOR THIS PATENT.

AB A method for treating hair loss in mammals uses compositions containing 2-decarboxy-2-phosphinico prostaglandin derivatives. The compositions can be applied topically to the skin. The compositions can arrest hair loss, reverse hair loss, and promote hair growth. Compositions containing 2-decarboxy-2-phosphinico prostaglandin derivatives can also be used to lower intraocular pressure and treat bone disorders.

CAS INDEXING IS AVAILABLE FOR THIS PATENT.

365241-18-5P 365241-19-6P 365241-20-9P 365241-21-0P 365241-22-1P 365241-23-2P 365241-24-3P 365241-25-4P 365241-26-5P 365241-27-6P 365241-28-7P 365241-29-8P 365241-30-1P 365241-31-2P 365241-32-3P 365241-33-4P 365241-34-5P 365241-35-6P 365241-36-7P 365241-37-8P 365241-38-9P 365241-39-0P 365241-40-3P 365241-41-4P 365241-45-8P 365241-43-6P 365241-47-0P 365241-45-8P 365241-49-2P 365241-50-5P 365241-51-6P 365241-55-0P 365241-50-1P 365241-57-2P 365241-58-3P 365241-59-4P 365241-60-7P

(cosmetic and pharmaceutical compns. contg. 2-decarboxy-2-phosphinico prostaglandin derivs.)

IT 365241-18-5P

(cosmetic and pharmaceutical compns. contg. 2-decarboxy-2-phosphinico prostaglandin derivs.)

RN 365241.-18-5 USPATFULL

CN Phosphinic acid, [6-[(1R,5S)-2-[(3R)-4-[(3-fluorophenyl)thio]-3-hydroxybutyl]-5-hydroxy-3-(hydroxyimino)cyclopentyl]hexyl]methyl- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Double bond geometry unknown.

=> fil reg FILE 'REGISTRY' ENTERED AT 07:11:06 ON 29 MAR 2002 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2002 American Chemical Society (ACS)

STRUCTURE FILE UPDATES: 27 MAR 2002 HIGHEST RN 403476-73-3 DICTIONARY FILE UPDATES: 27 MAR 2002 HIGHEST RN 403476-73-3

TSCA INFORMATION NOW CURRENT THROUGH July 7, 2001

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Calculated physical property data is now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details: http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf

The P indicator for Preparations was not generated for all of the CAS Registry Numbers that were added to the H/Z/CA/CAplus files between 12/27/01 and 1/23/02. Use of the P indicator in online and SDI searches during this period, either directly appended to a CAS Registry Number or by qualifying an L-number with /P, may have yielded incomplete results. As of 1/23/02, the situation has been resolved. Also, note that searches conducted using the PREP role indicator were not affected.

Customers running searches and/or SDIs in the H/Z/CA/CAplus files incorporating CAS Registry Numbers with the P indicator between 12/27/01 and 1/23/02, are encouraged to re-run these strategies. Contact the CAS Help Desk at 1-800-848-6533 in North America or 1-614-447-3698, worldwide, or send an e-mail to help@cas.org for further assistance or to receive a credit for any duplicate searches.

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=> s 121 or 123
L41 56 L21 OR L23
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=> d his 141-

(FILE 'HCAPLUS, USPATFULL' ENTERED AT 07:09:48 ON 29 MAR 2002)

FILE 'REGISTRY' ENTERED AT 07:10:08 ON 29 MAR 2002

FILE 'HCAPLUS, USPATFULL, USPAT2' ENTERED AT 07:10:30 ON 29 MAR 2002

FILE 'REGISTRY' ENTERED AT 07:11:06 ON 29 MAR 2002 L41 56 S L21 OR L23

FILE 'HCAPLUS' ENTERED AT 07:11:52 ON 29 MAR 2002 SEL HIT RN L34

FILE 'REGISTRY' ENTERED AT 07:11:57 ON 29 MAR 2002 64 S E1-E64

L42 64 S E1-E64 L43 8 S L42 NOT L41

=> d ide can tot 142

L42 ANSWER 1 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 380153-44-6 REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R,3R,5S)-3,5-bis(acetyloxy)-2-[(1E,3S)-3-(acetyloxy)-1-octenyl]cyclopentyl]-4-hexenyl]- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C25 H41 O8 P

SR CA

LC STN Files: CA, CAPLUS

Double bond geometry as shown.

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 136:31838

L42 ANSWER 2 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 380153-43-5 REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R,3R,5S)-3,5-bis(acetyloxy)-2-[(1E,3S)-3-(acetyloxy)-1-octenyl]cyclopentyl]-4-hexenyl]butyl-, ethyl ester (9CI) (CA INDEX NAME)

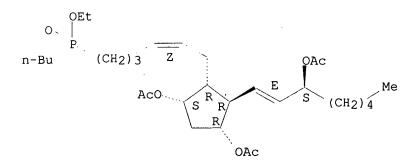
FS STEREOSEARCH

MF C31 H53 O8 P

SR CA

LC STN Files: CA, CAPLUS

Absolute stereochemistry. Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 136:31838

L42 ANSWER 3 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 380153-42-4 REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R,3R,5S)-3,5-bis(acetyloxy)-2-[(1E,3S)-3-(acetyloxy)-1-octenyl]cyclopentyl]-4-hexenyl]ethyl-, ethyl ester (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C29 H49 O8 P

SR CA

LC STN Files: CA, CAPLUS

Absolute stereochemistry. Double bond geometry as shown.

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 136:31838

L42 ANSWER 4 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 380153-41-3 REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R,3R,5S)-3,5-bis(acetyloxy)-2-[(1E,3S)-3-(acetyloxy)-1-octenyl]cyclopentyl]-4-hexenyl]methyl-, ethyl ester (9CI) (CA INDEX NAME)

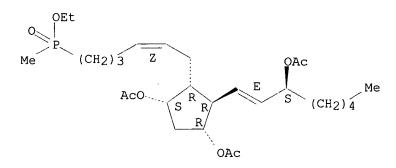
FS STEREOSEARCH

MF C28 H47 O8 P

SR CA

LC STN Files: CA, CAPLUS

Absolute stereochemistry. Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 136:31838

L42 ANSWER 5 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **380153-37-7** REGISTRY

CN Phosphinic acid, butyl[(4Z)-6-[(1R,2R,3R,5S)-3,5-dihydroxy-2-[(1E,3S)-3-hydroxy-1-octenyl]cyclopentyl]-4-hexenyl]- (9CI) (CA INDEX NAME)

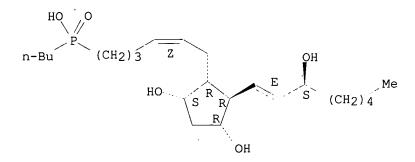
FS STEREOSEARCH

MF C23 H43 O5 P

SR CA

LC STN Files: CA, CAPLUS

Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 136:31838

L42 ANSWER 6 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 380153-36-6 REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R,3R,5S)-3,5-dihydroxy-2-[(1E,3S)-3-hydroxy-1-octenyl]cyclopentyl]-4-hexenyl]ethyl- (9CI) (CA INDEX NAME)

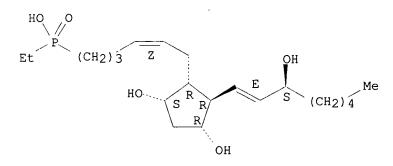
FS STEREOSEARCH

MF C21 H39 O5 P

SR CA

LC STN Files: CA, CAPLUS

Absolute stereochemistry. Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 136:31838

L42 ANSWER 7 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **380153-35-5** REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R,3R,5S)-3,5-dihydroxy-2-[(1E,3S)-3-hydroxy-1-octenyl]cyclopentyl]-4-hexenyl]- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C19 H35 O5 P

SR CA

LC STN Files: CA, CAPLUS

Double bond geometry as shown.

HO
$$\frac{P}{H}$$
 (CH2) $\frac{Z}{Z}$ OH $\frac{E}{R}$ OH OH

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 136:31838

L42 ANSWER 8 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 380153-34-4 REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R,3R,5S)-3,5-dihydroxy-2-[(1E,3S)-3-hydroxy-1-octenyl]cyclopentyl]-4-hexenyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C20 H37 O5 P

SR CA

LC STN Files: CA, CAPLUS

Absolute stereochemistry. Double bond geometry as shown.

HO O Me
$$P$$
 (CH2) $\frac{1}{3}$ $\frac{1}{Z}$ $\frac{1}{Z$

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 136:31838

L42 ANSWER 9 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **380153-33-3** REGISTRY

CN Phosphonic acid, [(4Z)-6-[(1R,2R,3R,5S)-3,5-dihydroxy-2-[(1E,3S)-3-hydroxy-1-octenyl]cyclopentyl]-4-hexenyl]- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C19 H35 O6 P

CI COM

SR CA

LC STN Files: CA, CAPLUS

Absolute stereochemistry.

Double bond geometry as shown.

$$H_2O_3P$$
 (CH₂) $\frac{1}{3}$ Z OH

 CH_2) $\frac{1}{4}$ R OH

1 REFERENCES IN FILE CA (1967 TO DATE) 1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

1: 136:31838 REFERENCE

ANSWER 10 OF 64 REGISTRY COPYRIGHT 2002 ACS L42

380153-32-2 REGISTRY RN

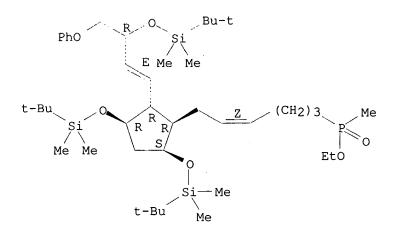
Phosphinic acid, [(4Z)-6-[(1R,2R,3R,5S)-3,5-bis[[(1,1-CN $\label{lem:dimethylethyl} \verb|dimethylsilyl| oxy| -2 - [(1E, 3R) -3 - [[(1, 1-1)]] -2 - [(1E, 3R) -3 - [[(1E, 3R)]] -2 - [(1E, 3R) -3 - [[(1E, 3R)]] -2 - [(1E, 3R)] -2 - [(1E$ dimethylethyl)dimethylsilyl]oxy}-4-phenoxy-1-butenyl]cyclopentyl]-4hexenyl]methyl-, ethyl ester (9CI) (CA INDEX NAME) FS STEREOSEARCH

C42 H79 O6 P Si3 MF

SR CA

LCSTN Files: CA, CAPLUS

Absolute stereochemistry. Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE) 1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 136:31838

ANSWER 11 OF 64 REGISTRY COPYRIGHT 2002 ACS L42

RN 380153-29-7 REGISTRY

Phosphinic acid, [(4Z)-6-[(1R,2R,3R,5S)-3,5-dihydroxy-2-[(1E,3R)-3-hydroxy-CN 4-phenoxy-1-butenyl]cyclopentyl]-4-hexenyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C22 H33 O6 P

SR CA

LC STN Files: CA, CAPLUS Absolute stereochemistry. Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 136:31838

L42 ANSWER 12 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-60-7 REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R,3R,5S)-3,5-dihydroxy-2-[(1E)-3-hydroxy-1-octenyl]cyclopentyl]-4-hexenyl]- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C19 H35 O5 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry. Double bond geometry as shown.

HO
$$\stackrel{P}{H}$$
 (CH₂) $\stackrel{Z}{3}$ $\stackrel{Z}{Z}$ OH $\stackrel{E}{K}$ (CH₂) $\stackrel{A}{4}$ Me

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 13 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **365241-59-4** REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R,3R,5S)-3,5-dihydroxy-2-[(1E)-3-hydroxy-4-phenoxy-1-butenyl]cyclopentyl]-4-hexenyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C22 H33 O6 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 14 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-58-3 REGISTRY

CN Phosphinic acid, [6-[(1R,2R,5S)-3,5-dihydroxy-2-[(3R)-3-hydroxy-4-(phenylamino)butyl]cyclopentyl]hexyl]- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C21 H36 N O5 P

SR - CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

1 REFERENCES IN FILE CA (1967 TO DATE)
1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 15 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-57-2 REGISTRY

CN Phosphinic acid, [6-[(1R,2R,5S)-3,5-dihydroxy-2-[(3R)-3-hydroxy-4-(phenylamino)butyl]cyclopentyl]hexyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C22 H38 N O5 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 16 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-56-1 REGISTRY

CN Phosphinic acid, [6-[(1R,2R,5S)-3,5-dihydroxy-2-[(3R)-3-hydroxy-5-phenylpentyl]cyclopentyl]hexyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C23 H39 O5 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 17 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-55-0 REGISTRY

CN Phosphinic acid, [6-[(1R,2R,5S)-3,5-dihydroxy-2-[(3R)-3-hydroxy-4-[3-(trifluoromethyl)phenoxy]butyl]cyclopentyl]hexyl]ethyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C24 H38 F3 O6 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 18 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-54-9 REGISTRY

CN Phosphinic acid, [6-[(1R,2R,5S)-2-[(3R)-4-(2-fluorophenoxy)-3-hydroxybutyl]-3,5-dihydroxycyclopentyl]hexyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C22 H36 F O6 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 19 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **365241-53-8** REGISTRY

CN Phosphinic acid, [6-[(1R,2R,5S)-3,5-dihydroxy-2-[(3R)-3-hydroxy-4-(phenylthio)butyl]cyclopentyl]hexyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C22 H37 O5 P S

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 20 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-52-7 REGISTRY

CN Phosphinic acid, [6-[(1R,2R,5S)-3,5-dihydroxy-2-[(3R)-3-hydroxy-6-phenyl-4,5-hexadienyl]cyclopentyl]hexyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C24 H37 O5 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

$$Ph$$
 $C = C = C$ R OH P OH OH OH

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 21 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **365241-51-6** REGISTRY

CN Phosphinic acid, [6-[(1R,2R,5S)-3,5-dihydroxy-2-[(3R)-3-hydroxy-5-phenyl-4-pentynyl]cyclopentyl]hexyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C23 H35 O5 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 22 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-50-5 REGISTRY

CN Phosphinic acid, [6-[(1R,2R,5S)-3,5-dihydroxy-2-[(3R)-3-hydroxy-5-phenyl-4-pentenyl]cyclopentyl]hexyl]ethyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C24 H39 O5 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

Double bond geometry unknown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 23 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **365241-49-2** REGISTRY

CN Phosphinic acid, [6-[(1R,2R,5S)-2-[(3R)-4-[(2-fluorophenyl)thio]-3-hydroxybutyl]-3,5-dihydroxycyclopentyl]hexyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C22 H36 F O5 P S

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 24 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **365241-48-1** REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R)-2-[(1E,3R)-4-(3-chlorophenoxy)-3-hydroxy-1-butenyl]-3,5-dihydroxycyclopentyl]-4-hexenyl]ethyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C23 H34 C1 O6 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry. Double bond geometry as shown.

OH
$$(CH_2)_3$$
 P Et $(CH_2)_4$ HO

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 25 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-47-0 REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R)-3,5-dihydroxy-2-[(1E,3S)-3-hydroxy-4-phenyl-1-butenyl]cyclopentyl]-4-hexenyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C22 H33 O5 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 26 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-46-9 REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R)-3,5-dihydroxy-2-[(1E,3S)-3-hydroxy-5-phenyl-1-pentenyl]cyclopentyl]-4-hexenyl]propyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C25 H39 O5 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry. Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 27 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **365241-45-8** REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R)-3,5-dihydroxy-2-[(1E,3R)-3-hydroxy-4-[3-(trifluoromethyl)phenoxy]-1-butenyl]cyclopentyl]-4-hexenyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C23 H32 F3 O6 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry. Double bond geometry as shown.

F3C OH
$$(CH_2)_3$$
 P Me

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 28 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **365241-44-7** REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R)-3,5-dihydroxy-2-[(1E,3S)-3-hydroxy-1,5-octadienyl]cyclopentyl]-4-hexenyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C20 H35 O5 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

Double bond geometry as described by E or Z.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 29 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **365241-43-6** REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R)-3,5-dihydroxy-2-[(1E,3S)-3-hydroxy-1-octenyl]cyclopentyl]-4-hexenyl]ethyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C21 H39 O5 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 30 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-42-5 REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R)-3,5-dihydroxy-2-[(1E,3S)-3-hydroxy-1-octenyl]cyclopentyl]-4-hexenyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C20 H37 O5 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry. Double bond geometry as shown.

Me
$$P$$
 (CH_2) $\frac{Z}{Z}$ P OH E S (CH_2) $\frac{A}{4}$ Me OH

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 31 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-41-4 REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R)-2-[(1E,3R)-4-[(2,4-difluorophenyl)amino]-3-hydroxy-1-butenyl]-3,5-dihydroxycyclopentyl]-4-hexenyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C22 H32 F2 N O5 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry. Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 32 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-40-3 REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R)-2-[(1E,3R)-4-[(2-fluorophenyl)amino]-3-hydroxy-1-butenyl]-3,5-dihydroxycyclopentyl]-4-hexenyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C22 H33 F N O5 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)
1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 33 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-39-0 REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R)-2-[(1E,3S)-5-[1,1'-biphenyl]-4-yl-3-hydroxy-1-pentenyl]-3,5-dihydroxycyclopentyl]-4-hexenyl]propyl- (9CI) (CA

INDEX NAME)

FS STEREOSEARCH

MF C31 H43 O5 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

Double bond geometry as shown.

OH
$$CH_{2})_{3}$$

$$Pr-n$$

$$R$$

$$R$$

$$OH$$

$$R$$

$$R$$

$$OH$$

$$R$$

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 34 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **365241-38-9** REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R)-3,5-dihydroxy-2-[(1E,3R)-3-hydroxy-4-[3-(trifluoromethyl)phenoxy]-1-butenyl]cyclopentyl]-4-hexenyl]ethyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C24 H34 F3 O6 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

Double bond geometry as shown.

F3C OH
$$R$$
 R OH R OH R R OH

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 35 OF 64 REGISTRY COPYRIGHT 2002 ACS RN 365241-37-8 REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R)-3,5-dihydroxy-2-[(1E,3R)-3-hydroxy-4-phenoxy-1-butenyl]cyclopentyl]-4-hexenyl]ethyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C23 H35 O6 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 36 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-36-7 REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R)-2-[(1E,3R)-4-(3-chlorophenoxy)-3-hydroxy-1-butenyl]-3,5-dihydroxycyclopentyl]-4-hexenyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C22 H32 C1 O6 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 37 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-35-6 REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R)-2-[(1E,3R)-4-[(2-fluorophenyl)thio]-3-hydroxy-1-butenyl]-3,5-dihydroxycyclopentyl]-4-hexenyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C22 H32 F O5 P S

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 38 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-34-5 REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R)-2-[(1E,3S)-5-(2-fluorophenyl)-3-hydroxy-1-pentenyl]-3,5-dihydroxycyclopentyl]-4-hexenyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C23 H34 F O5 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

- 1 REFERENCES IN FILE CA (1967 TO DATE)
- 1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 39 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-33-4 REGISTRY

CN Phosphinic acid, [6-[(1R,2R,5S)-2-[4-[(2-fluorophenyl)thio]-3-hydroxybutyl]-5-hydroxy-3-(hydroxyamino)cyclopentyl]hexyl](1-methylethyl)-(9CI) (CA INDEX NAME)

STEREOSEARCH

MF C24 H41 F N O5 P S

SR CA

FS

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 40 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **365241-32-3** REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R,5S)-2-[(1E)-4-[(4,5-dichloro-2-thienyl)thio]-3-hydroxy-1-butenyl]-5-hydroxy-3-(hydroxyamino)cyclopentyl]-4-hexenyl]ethyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C21 H32 C12 N O5 P S2

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 41 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-31-2 REGISTRY

CN Phosphinic acid, ethyl[6-[(1R,2R,5S)-5-hydroxy-3-(hydroxyamino)-2-[3-hydroxy-4-(phenylamino)butyl]cyclopentyl]hexyl]- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C23 H41 N2 O5 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 42 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **365241-30-1** REGISTRY

CN Phosphinic acid, [6-[(1R,2R,5S)-2-(3-benzo[b]thien-2-yl-3-hydroxypropyl)-5-hydroxy-3-(hydroxyamino)cyclopentyl]hexyl]ethyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C24 H38 N O5 P S

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 43 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **365241-29-8** REGISTRY

CN Phosphinic acid, [6-[(1R,2R,5S)-2-[5-(2-fluorophenyl)-3-hydroxy-4-pentynyl]-5-hydroxy-3-(hydroxyamino)cyclopentyl]hexyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C23 H35 F N O5 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 44 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **365241-28-7** REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R,5S)-5-hydroxy-3-(hydroxyamino)-2-[(1E)-3-hydroxy-3-phenyl-1-propenyl]cyclopentyl]-4-hexenyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C21 H32 N O5 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

- 1 REFERENCES IN FILE CA (1967 TO DATE)
- 1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 45 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-27-6 REGISTRY

CN Phosphinic acid, butyl[6-[(1R,5S)-2-[(3R)-5-(2-fluorophenyl)-3-hydroxy-4-pentynyl]-5-hydroxy-3-(hydroxyimino)cyclopentyl]hexyl]- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C26 H39 F N O5 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry. Double bond geometry unknown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 46 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **365241-26-5** REGISTRY

CN Phosphinic acid, [6-[(1R,5S)-2-[(3R)-4-(benzo[b]thien-3-ylthio)-3-hydroxybutyl]-5-hydroxy-3-(hydroxyimino)cyclopentyl]hexyl]ethyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C25 H38 N O5 P S2

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

Double bond geometry unknown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

- 1 REFERENCES IN FILE CA (1967 TO DATE)
- 1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 47 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-25-4 REGISTRY

CN Phosphinic acid, [6-[(1R,5S)-5-hydroxy-2-[(3S)-3-hydroxyoctyl]-3-(methoxyimino)cyclopentyl]hexyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C21 H42 N O5 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

Double bond geometry unknown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 48 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **365241-24-3** REGISTRY

CN Phosphinic acid, [6-[(1R,5S)-2-[(3R)-4-[(2,4-difluorophenyl)amino]-3-hydroxybutyl]-5-hydroxy-3-(hydroxyimino)cyclopentyl]hexyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C22 H35 F2 N2 O5 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry. Double bond geometry unknown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 49 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-23-2 REGISTRY

CN Phosphinic acid, [6-[(1R,5S)-2-[(3R)-4-(3-chlorophenoxy)-3-hydroxybutyl]-5-hydroxy-3-(hydroxyimino)cyclopentyl]hexyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C22 H35 C1 N O6 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry. Double bond geometry unknown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 50 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **365241-22-1** REGISTRY

CN Phosphinic acid, [6-[(1R,5S)-5-hydroxy-3-(hydroxyimino)-2-[(3S)-3-hydroxynonyl]cyclopentyl]hexyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C21 H42.N O5 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.
Double bond geometry unknown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 51 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-21-0 REGISTRY

CN Phosphinic acid, [6-[(1R,5S)-5-hydroxy-3-(hydroxyimino)-2-[(3S)-3-hydroxyoctyl]cyclopentyl]hexyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C20 H40 N O5 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry. Double bond geometry unknown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 52 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 365241-20-9 REGISTRY

CN Phosphinic acid, [6-[(1R,5S)-2-[(3R)-4-(3-fluorophenoxy)-3-hydroxybutyl]-5-hydroxy-3-(hydroxyimino)cyclopentyl]hexyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C22 H35 F N O6 P

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

Double bond geometry unknown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 53 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **365241-19-6** REGISTRY

CN Phosphinic acid, ethyl[6-[(1R,5S)-2-[(3R)-4-[(2-fluorophenyl)thio]-3-

hydroxybutyl]-5-hydroxy-3-(hydroxyimino)cyclopentyl]hexyl]- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C23 H37 F N O5 P S

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

Double bond geometry unknown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 54 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **365241-18-5** REGISTRY

CN Phosphinic acid, [6-[(1R,5S)-2-[(3R)-4-[(3-fluorophenyl)thio]-3-hydroxybutyl]-5-hydroxy-3-(hydroxyimino)cyclopentyl]hexyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C22 H35 F N O5 P S

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

Double bond geometry unknown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 135:293970

L42 ANSWER 55 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **325153-12-6** REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R, 2R, 3R, 5S)-3, 5-dihydroxy-2-[(1E, 3R)-3-hydroxy-

4-(2-methoxyethoxy)-1-butenyl]cyclopentyl]-4-hexenyl]methyl- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C19 H35 O7 P

SR CA

LC STN Files: CA, CAPLUS

Absolute stereochemistry. Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 134:162868

L42 ANSWER 56 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 325153-11-5 REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R,3R,5S)-2-[(1E,3S)-5-(2-fluorophenyl)-3-hydroxy-1-pentenyl]-3,5-dihydroxycyclopentyl]-4-hexenyl]methyl-, monosodium salt (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C23 H34 F O5 P . Na

SR CA

LC STN Files: CA, CAPLUS

Absolute stereochemistry. Double bond geometry as shown.

Na

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 134:162868

L42 ANSWER 57 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 252858-17-6 REGISTRY

CN Phosphonic acid, [(4Z)-6-[(1R,2R,3R,5S)-3,5-dihydroxy-2-[(1E,3S)-3-hydroxy-1-octenyl]cyclopentyl]-4-hexenyl]-, disodium salt (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C19 H35 O6 P . 2 Na

SR CF

LC STN Files: CA, CAPLUS, CASREACT

CRN (380153-33-3)

Absolute stereochemistry. Double bond geometry as shown.

$$(CH_2)_3$$
 Z OH E S R $CH_2)_4$ Me OH

●2 Na

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 132:49813

L42 ANSWER 58 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **252858-16-5** REGISTRY

CN Phosphonic acid, [(4Z)-6-[(1R,2R,3R,5S)-3,5-dihydroxy-2-[(1E,3S)-3-hydroxy-1-octenyl]cyclopentyl]-4-hexenyl]-, monoethyl ester (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C21 H39 O6 P

SR CA

LC STN Files: CA, CAPLUS, CASREACT

Absolute stereochemistry.

Double bond geometry as shown.

Eto P (CH₂) $\frac{Z}{Z}$ OH $\frac{E}{R}$ OH OH

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

2 REFERENCES IN FILE CA (1967 TO DATE)

2 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 136:31838

REFERENCE 2: 132:49813

L42 ANSWER 59 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 252858-14-3 REGISTRY

CN Phosphonic acid, [(4Z)-6-[(1R,2R,3R,5S)-3,5-bis(acetyloxy)-2-[(1E,3S)-3-(acetyloxy)-1-octenyl]cyclopentyl]-4-hexenyl]-, diethyl ester (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C29 H49 O9 P

SR CA

LC STN Files: CA, CAPLUS, CASREACT

Absolute stereochemistry. Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 132:49813

L42 ANSWER 60 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **252858-12-1** REGISTRY

CN Phosphonic acid, [6-[(1R,2R,3R,5S)-3,5-dihydroxy-2-[(1E,3S)-3-hydroxy-1-octenyl]cyclopentyl]hexyl]-, monoethyl ester (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C21 H41 O6 P

SR CA

LC STN Files: CA, CAPLUS, CASREACT

Absolute stereochemistry.

Double bond geometry as shown.

HO O
$$(CH_2)_6$$
 OH E S R R $CH_2)_4$ Me

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 132:49813

L42 ANSWER 61 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 252858-11-0 REGISTRY

CN Phosphonic acid, [6-[(1R,2R,3R,5S)-3,5-bis(acetyloxy)-2-[(1E,3S)-3-(acetyloxy)-1-octenyl]cyclopentyl]hexyl]-, diethyl ester (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C29 H51 O9 P

SR CA

LC STN Files: CA, CAPLUS, CASREACT

Absolute stereochemistry. Double bond geometry as shown.

OEt
OP
EtO
(CH2)6
OAC

ACO
SR
R
OAC
OAC

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 132:49813

L42 ANSWER 62 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN **77493-13-1** REGISTRY

CN Phosphonic acid, [(9.alpha.,11.alpha.,13E,15S)-9,11,15-tris[(tetrahydro-2H-pyran-2-yl)oxy]prost-13-en-1-yl]-, dimethyl ester (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:

CN Prostane, phosphonic acid deriv.

FS STEREOSEARCH

MF C37 H67 O9 P

LC STN Files: CA, CAPLUS

Absolute stereochemistry.

Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 94:191740

L42 ANSWER 63 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 72489-45-3 REGISTRY

CN Phosphonic acid, [(9.beta., 11.alpha., 13E, 15S)-9, 11, 15-trihydroxyprost-13-

en-1-yl]-, dimethyl ester (9CI) (CA INDEX NAME)

OTHER CA INDEX NAMES:

CN Prostane, phosphonic acid deriv.

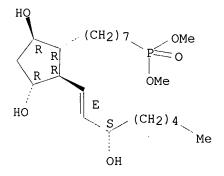
FS STEREOSEARCH

MF C22 H43 O6 P

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

2 REFERENCES IN FILE CA (1967 TO DATE)

2 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 94:191740

REFERENCE 2: 92:58302

L42 ANSWER 64 OF 64 REGISTRY COPYRIGHT 2002 ACS

RN 72488-21-2 REGISTRY

CN Phosphonic acid, [(9.alpha.,11.alpha.,13E,15S)-9,11,15-trihydroxyprost-13-en-1-yl]-, dimethyl ester (9CI) (CA INDEX NAME)

OTHER CA INDEX NAMES:

CN Prostane, phosphonic acid deriv.

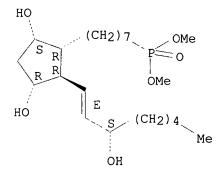
FS STEREOSEARCH

MF C22 H43 O6 P

LC STN Files: CA, CAPLUS, USPATFULL

Absolute stereochemistry.

Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

2 REFERENCES IN FILE CA (1967 TO DATE)

2 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 94:191740

REFERENCE 2: 92:58302

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L43 ANSWER 1 OF 8 REGISTRY COPYRIGHT 2002 ACS

RN 380153-44-6 REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R,3R,5S)-3,5-bis(acetyloxy)-2-[(1E,3S)-3-(acetyloxy)-1-octenyl]cyclopentyl]-4-hexenyl]- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

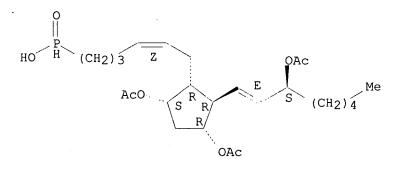
MF C25 H41 O8 P

SR CA

LC STN Files: CA, CAPLUS

Absolute stereochemistry.

Double bond geometry as shown.



- 1 REFERENCES IN FILE CA (1967 TO DATE)
- 1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 136:31838

L43 ANSWER 2 OF 8 REGISTRY COPYRIGHT 2002 ACS

RN 380153-43-5 REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R,3R,5S)-3,5-bis(acetyloxy)-2-[(1E,3S)-3-(acetyloxy)-1-octenyl]cyclopentyl]-4-hexenyl]butyl-, ethyl ester (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C31 H53 O8 P

SR CA

LC STN Files: CA, CAPLUS

Absolute stereochemistry. Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 136:31838

L43 ANSWER 3 OF 8 REGISTRY COPYRIGHT 2002 ACS

RN 380153-42-4 REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R,3R,5S)-3,5-bis(acetyloxy)-2-[(1E,3S)-3-(acetyloxy)-1-octenyl]cyclopentyl]-4-hexenyl]ethyl-, ethyl ester (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C29 H49 O8 P

SR CA

LC STN Files: CA, CAPLUS

Absolute stereochemistry.

Double bond geometry as shown.

1 REFERENCES IN FILE CA (1967 TO DATE)
1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 136:31838

L43 ANSWER 4 OF 8 REGISTRY COPYRIGHT 2002 ACS

RN 380153-41-3 REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R,3R,5S)-3,5-bis(acetyloxy)-2-[(1E,3S)-3-(acetyloxy)-1-octenyl]cyclopentyl]-4-hexenyl]methyl-, ethyl ester (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C28 H47 O8 P

SR CA

LC STN Files: CA, CAPLUS

Absolute stereochemistry. Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 136:31838

L43 ANSWER 5 OF 8 REGISTRY COPYRIGHT 2002 ACS

RN **380153-32-2** REGISTRY

CN Phosphinic acid, [(4Z)-6-[(1R,2R,3R,5S)-3,5-bis[[(1,1-dimethylethyl)dimethylsilyl]oxy]-2-[(1E,3R)-3-[[(1,1-dimethylethyl)dimethylsilyl]oxy]-4-phenoxy-1-butenyl]cyclopentyl]-4-hexenyl]methyl-, ethyl ester (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C42 H79 O6 P Si3

SR CA

LC STN Files: CA, CAPLUS

Absolute stereochemistry. Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 136:31838

L43 ANSWER 6 OF 8 REGISTRY COPYRIGHT 2002 ACS

RN 252858-14-3 REGISTRY

CN Phosphonic acid, [(4Z)-6-[(1R,2R,3R,5S)-3,5-bis(acetyloxy)-2-[(1E,3S)-3-(acetyloxy)-1-octenyl]cyclopentyl]-4-hexenyl]-, diethyl ester (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C29 H49 O9 P

SR CA

LC STN Files: CA, CAPLUS, CASREACT

Absolute stereochemistry. Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 132:49813

L43 ANSWER 7 OF 8 REGISTRY COPYRIGHT 2002 ACS

RN 252858-11-0 REGISTRY

CN Phosphonic acid, [6-[(1R,2R,3R,5S)-3,5-bis(acetyloxy)-2-[(1E,3S)-3-(acetyloxy)-1-octenyl]cyclopentyl]hexyl]-, diethyl ester (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C29 H51 O9 P

SR CF

LC STN Files: CA, CAPLUS, CASREACT

Absolute stereochemistry. Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 132:49813

L43 ANSWER 8 OF 8 REGISTRY COPYRIGHT 2002 ACS

RN **77493-13-1** REGISTRY

CN Phosphonic acid, [(9.alpha.,11.alpha.,13E,15S)-9,11,15-tris[(tetrahydro-2H-pyran-2-yl)oxy]prost-13-en-1-yl]-, dimethyl ester (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:

CN Prostane, phosphonic acid deriv.

FS STEREOSEARCH

MF C37 H67 O9 P

LC STN Files: CA, CAPLUS

Absolute stereochemistry. Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

- 1 REFERENCES IN FILE CA (1967 TO DATE)
- 1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 94:191740